Felice Grandinetti

List of Publications by Year
in descending order

121	1,989	257450	361022
papers	citations	24 h-index	35 g-index

127	127	127	841
all docs	docs citations	times ranked	citing authors

#	Article	IF	Citations
1	Noble-gas compounds: A general procedure of bonding analysis. Journal of Chemical Physics, 2022, 156, 014104.	3.0	10
2	On the Proton-Bound Noble Gas Dimers (Ng-H-Ng)+ and (Ng-H-Ngâ€™)+ (Ng, Ngâ€™ = He-Xe): Relationships between Structure, Stability, and Bonding Character. Molecules, 2021, 26, 1305.	3.8	8
3	From LAr to L-ArBeO (LÂ=ÂHe, Ne, Ar, HF): Switching on Ïƒ-hole effects in non-covalent interactions. Chemical Physics Letters, 2021, 768, 138402.	2.6	8
4	Concerning the Role of Ïƒ-Hole in Non-Covalent Interactions: Insights from the Study of the Complexes of ArBeO with Simple Ligands. Molecules, 2021, 26, 4477.	3.8	2
5	Classifying the chemical bonds involving the noble-gas atoms. New Journal of Chemistry, 2020, 44, 14536-14550.	2.8	17
6	Cationic Noble-Gas Hydrides: From Ion Sources to Outer Space. Frontiers in Chemistry, 2020, 8, 462.	3.6	21
7	Complexes of helium with neutral molecules: Progress toward a quantitative scale of bonding character. Journal of Computational Chemistry, 2020, 41, 1000-1011.	3.3	10
8	Complexes of the noble-gas atoms with unsaturated ions: A theoretical investigation on the exemplary (H2CÂ=ÂNH2+)Ar. Chemical Physics Letters, 2020, 752, 137532.	2.6	1
9	Noncovalent Complexes of the Nobleâ€Gas Atoms: Analyzing the Transition from Physical to Chemical Interactions. Journal of Computational Chemistry, 2019, 40, 2318-2328.	3.3	19
10	Helium Accepts Back-Donation In Highly Polar Complexes: New Insights into the Weak Chemical Bond. Journal of Physical Chemistry Letters, 2017, 8, 3334-3340.	4.6	24
11	Complexes of the Noble Gases with H3O+: A Theoretical Investigation of Ng(H3O+) (Ng = Heâ€”Xe). European Journal of Mass Spectrometry, 2015, 21, 171-181.	1.0	7
12	Bimolecular Homolytic Substitutions at Nitrogen: An Experimental and Theoretical Study on the Gasâ€Phase Reactions of Alkyl Radicals with NF₃. Chemistry - A European Journal, 2015, 21, 15826-15834.	3.3	5
13	Catching the role of anisotropic electronic distribution and charge transfer in halogen bonded complexes of noble gases. Journal of Chemical Physics, 2015, 142, 184304.	3.0	39
14	Bonding Motifs of Noble-Gas Compounds As Described by the Local Electron Energy Density. Journal of Physical Chemistry A, 2015, 119, 6528-6541.	2.5	42
15	Electronic structure and conformational flexibility of d-cycloserine. Physical Chemistry Chemical Physics, 2015, 17, 25845-25853.	2.8	1
16	Experimental Evidence of Chemical Components in the Bonding of Helium and Neon with Neutral Molecules. Chemistry - A European Journal, 2015, 21, 6234-6240.	3.3	53
17	Complexes of XeHXe⁺ with Simple Ligands: A Theoretical Investigation on (XeHXe⁺)L (L = N₂, CO, H₂O, NH₃). Journal of Physical Chemistry A, 2015, 119, 2383-2392.	2.5	22
18	Neutral Compounds with Xenonâ€“Germanium Bonds: A Theoretical Investigation on FXeGeF and FXeGeF3. Journal of Physical Chemistry A, 2014, 118, 3326-3334.	2.5	21

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19	Ion chemistry of sulfuryl fluoride: An experimental and theoretical study on gas-phase reactions involving neutral and ionized SO2F2. International Journal of Mass Spectrometry, 2013, 354-355, 46-53.	1.5	0
20	Neon behind the signs. Nature Chemistry, 2013, 5, 438-438.	13.6	25
21	<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" altimg="si1.gif" overflow="scroll"><mml:mrow><mml:mo stretchy="false">(</mml:mo><mml:msup><mml:mrow><mml:mtext>HNg</mml:mtext></mml:mrow><mml:mrow><mml:mo>+</mml:mo stretchy="false">(</mml:mo><mml:msub><mml:mrow><mml:mtext>OH</mml:mtext></mml:mrow><mml:mrow><mml:mn>2</mml:mn	2.5	5
22	Germyl Cations with Geâ€“S Bonds: An Experimental and Theoretical Study on the Gaseous F_nGe(SH)⁺_{3â€“n} (n = 0â€“2). European Journal of Mass Spectrometry, 2012, 18, 447-456.	1.0	1
23	Gaseous germyl cations: A theoretical investigation on the structure, properties, and mechanism of formation of and (n=0â€“2). Computational and Theoretical Chemistry, 2012, 993, 131-139.	2.5	6
24	Gasâ€phase reactions of SiH_n⁺ (n = 1,2) with NF₃: A computational investigation on the detailed mechanistic aspects. Journal of Computational Chemistry, 2012, 33, 1918-1926.	3.3	3
25	Gas-Phase Ion Chemistry of the Noble Gases: Recent Advances and Future Perspectives. European Journal of Mass Spectrometry, 2011, 17, 423-463.	1.0	69
26	Positive Ion Chemistry of SiH₄/GeF₄ Gaseous Mixtures Studied by Ion Trap Mass Spectrometry and Ab Initio Calculations. European Journal of Mass Spectrometry, 2011, 17, 197-206.	1.0	2
27	Gasâ€phase chemistry of ionized and protonated GeF₄: a joint experimental and theoretical study. Journal of Mass Spectrometry, 2011, 46, 465-477.	1.6	10
28	Xenonâ€“Nitrogen Chemistry: Gasâ€Phase Generation and Theoretical Investigation of the Xenonâ€“Difluoronitrenium Ion F₂Nï£¿Xe⁺. Chemistry - A European Journal, 2011, 17, 10682-10689.	3.3	40
29	Cationic noble gas hydrides-2: A theoretical investigation on HNgHNgH+ (Ng=Ar, Kr, Xe). Computational and Theoretical Chemistry, 2011, 964, 318-323.	2.5	19
30	Stabilization of HHeF by Complexation: Is it a Really Viable Strategy?. Chemistry - A European Journal, 2010, 16, 6257-6264.	3.3	7
31	F3Geâ”Xe+: A Xenonâ”Germanium Molecular Species. Journal of Physical Chemistry Letters, 2010, 1, 2006-2010.	4.6	39
32	Cationic Noble Gas Hydrides: A Theoretical Investigation of Dinuclear HNgFNgH⁺(Ng =) Tj ETQq0 0 0 rgBT /Overlock 10 Tf 5	2.5	29
33	Ion/Molecule reactions in SiH₄/H₂S and GeH₄/H₂S mixtures. Journal of Mass Spectrometry, 2009, 44, 725-734.	1.6	5
34	Gasâ€phase reactions of XH₃⁺ (X = C, Si, Ge) with NF₃: a comparative investigation on the detailed mechanistic aspects. Journal of Mass Spectrometry, 2009, 44, 1348-1358.	1.6	8
35	Protonated MF3 (M=Nâ€“Bi): Structure, stability, and thermochemistry of the Hâ€“MF3+ and HFâ€“MF2+ isomers. Journal of Fluorine Chemistry, 2009, 130, 557-561.	1.7	7
36	Noble gasâ€“selenium molecular species: A theoretical investigation of FNgSeâ” (Ng=Heâ€“Xe). Chemical Physics Letters, 2009, 470, 49-53.	2.6	18

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37	Positive Ion Chemistry of SiH4/NF3 Gaseous Mixtures Studied by Ion Trap Mass Spectrometry. European Journal of Mass Spectrometry, 2009, 15, 209-220.	1.0	7
38	Cl-Initiated oxidation of N-ethyl-perfluoroalkanesulfonamides: A theoretical insight into the experimentally observed products. Computational and Theoretical Chemistry, 2008, 857, 57-65.	1.5	1
39	Ion chemistry in germane/fluorocompounds gaseous mixtures: a mass spectrometric and theoretical study. Journal of Mass Spectrometry, 2008, 43, 1320-1333.	1.6	11
40	Noble gasâ€“sulfur anions: A theoretical investigation of FNgSâ¯ (Ng=He, Ar, Kr, Xe). Chemical Physics Letters, 2008, 458, 48-53.	2.6	29
41	Noble Gas Anions:â€‰ A Theoretical Investigation of FNgBN⁻(Ng = Heâ¯Xe). Journal of Physical Chemistry A, 2007, 111, 10144-10151.	2.5	53
42	Chemically enhanced liquid chromatography/tandem mass spectrometry determination of glutamic acid in the diffusion medium of retinal cells. Biomedical Chromatography, 2007, 21, 1069-1076.	1.7	29
43	Nobleâ€Gas Complexes: Theoretical Investigation of Multicenter Polynuclear Species. Helvetica Chimica Acta, 2007, 90, 1335-1352.	1.6	4
44	Gas-phase ion chemistry of NF3/SO2 mixtures: A mass spectrometric and theoretical investigation. International Journal of Mass Spectrometry, 2007, 266, 86-91.	1.5	2
45	Ge3Hn-Anions (n= 0â¯5) and Their Neutral Analogues:Â A Theoretical Investigation on the Structure, Stability, and Thermochemistry. Journal of Physical Chemistry A, 2006, 110, 9429-9437.	2.5	3
46	Cationic Germanium Fluorides:Â A Theoretical Investigation on the Structure, Stability, and Thermochemistry of GeFn/GeFn+(n= 1â¯3). Journal of Physical Chemistry A, 2006, 110, 4900-4905.	2.5	15
47	Ligation of Be+ and Mg+ to NF3: Structure, stability, and thermochemistry of the Be+â€“(NF3) and Mg+â€“(NF3) complexes. International Journal of Mass Spectrometry, 2006, 255-256, 11-19.	1.5	4
48	Cationic germanium fluorides. International Journal of Mass Spectrometry, 2006, 257, 50-59.	1.5	8
49	Neutral Helium Compounds: Theoretical Evidence for a Large Class of Polynuclear Complexes. Chemistry - A European Journal, 2006, 12, 5033-5042.	3.3	36
50	Fluoromethyl Cations and Group XIV Congeners AHnF3 â€“n+ (A = Si, Ge, Sn, Pb;n = 0â€“2): From Covalent Structures to Ion-Molecule Complexes. European Journal of Inorganic Chemistry, 2006, 2006, 3010-3015.	2.0	5
51	From OBeHe to H3BOBeHe: Enhancing the stability of a neutral helium compound. Chemical Physics Letters, 2005, 406, 179-183.	2.6	43
52	Helium Chemistry: A Survey of the Role of the Ionic Species. ChemInform, 2005, 36, no.	0.0	0
53	Comment on â€œComputational Investigation of SO3â¯NH3-nXn(n= 0â¯3; X = F, Cl) Interactionsâ€• Journal of Physical Chemistry A, 2005, 109, 2410-2411.	2.5	1
54	Nitrogen Trifluoride as a Bifunctional Lewis Base: Implications for the Adsorption of NF3 on Solid Surfaces. European Journal of Inorganic Chemistry, 2004, 2004, 1125-1130.	2.0	11

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55	A Computational Investigation of HCN2+ Isomeric Structures: Implications for the Chemistry of Titan's Atmosphere. ChemPhysChem, 2004, 5, 1345-1351.	2.1	3
56	FN+Cl Ions from Ionized F2NCl: a Computational Investigation on the Structure and Reactivity toward H2O. Helvetica Chimica Acta, 2004, 87, 1467-1482.	1.6	3
57	Helium chemistry: a survey of the role of the ionic species. International Journal of Mass Spectrometry, 2004, 237, 243-267.	1.5	93
58	SBeNg, SBNg+, and SCNg2+ complexes (Ng=He, Ne, Ar): a computational investigation on the structure and stability. Chemical Physics Letters, 2004, 384, 25-29.	2.6	38
59	OBHe+: a remarkably stable singly charged cation containing helium. Chemical Physics Letters, 2004, 398, 357-360.	2.6	6
60	Concerning the reaction between singlet nitrenium ions and water: A computational investigation on competitive reaction paths. Journal of Computational Chemistry, 2003, 24, 547-564.	3.3	5
61	Isomeric Alkyl Cation/Arene Complexes in the Gas Phase. Chemistry - A European Journal, 2003, 9, 2072-2078.	3.3	17
62	FSO+ and FSO2+ ions from ionised sulfur oxyfluorides: a computational investigation on the structure, stability, and thermochemistry. Chemical Physics Letters, 2003, 372, 455-463.	2.6	7
63	A computational investigation on the Lewis acidity of fluoro- and chloronitrenium ions. Computational and Theoretical Chemistry, 2003, 635, 221-227.	1.5	5
64	Berylliumâ€“helium cations: computational evidence for a large class of thermodynamically stable species. International Journal of Mass Spectrometry, 2003, 228, 415-427.	1.5	19
65	Stable Compounds of the Lightest Noble Gases:â€‰ A Computational Investigation of RNBeNg (Ng = He, Ne,) Tj ETQq1 1 0.784314 rgB	2.5	57
66	Adducts of NF2+ with diatomic and simple polyatomic ligands: a computational investigation on the structure, stability, and thermochemistry. International Journal of Mass Spectrometry, 2002, 216, 285-299.	1.5	24
67	A computational investigation on the mechanism of the reaction between O(1D) and NF3. Chemical Physics Letters, 2002, 366, 676-682.	2.6	9
68	Complexes of lithium cation with nitrogen trifluoride: a computational investigation on the structure and stability of Li+â€“(NF3) isomers. Computational and Theoretical Chemistry, 2001, 574, 185-193.	1.5	14
69	The Unimolecular Loss of HF by Simple Inorganic Ions: A Computational Dynamic Reaction Path Study. European Journal of Mass Spectrometry, 2000, 6, 31-37.	1.0	5
70	Eliminative Ring Opening of Oxiranium Ions in the Gas Phase. Angewandte Chemie - International Edition, 2000, 39, 1673-1676.	13.8	6
71	FBeNg+ (Ng=He, Ne, Ar): Suitable Cations for Salts of the Lightest Noble Gases?. Angewandte Chemie - International Edition, 2000, 39, 1690-1692.	13.8	12
72	Spin-forbidden F+ transfer between 2NF+ and CO: a computational study on the detailed mechanistic aspects. International Journal of Mass Spectrometry, 2000, 201, 151-160.	1.5	3

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73	The ionization potential of NF 3 : a G3 computational study on the thermochemical properties of NF x and NF x + (x =1â€“3). Computational and Theoretical Chemistry, 2000, 497, 205-209.	1.5	16
74	Unimolecular decay of the thiomethoxy cation, CH3S+: A computational study on the detailed mechanistic aspects. Journal of Chemical Physics, 1999, 111, 6759-6768.	3.0	19
75	Carbonylation of ammonia by gaseous FCO+. A G2 and Rice-Ramsperger-Kassel-Marcus study of the detailed mechanistic aspects. International Journal of Mass Spectrometry, 1999, 184, 89-101.	1.5	5
76	Mechanistic Aspects of F+ Transfer Reactions: A Model Study in the Gas Phase. Chemistry - A European Journal, 1998, 4, 2366-2374.	3.3	19
77	Protonated NF3O. A G2MS theoretical study on the structure, stability, and interconversion of the (NF3O)H+ isomers. International Journal of Mass Spectrometry and Ion Processes, 1998, 175, 317-324.	1.8	4
78	Chiral Ions in the Gas Phase. 1. Intramolecular Racemization and Isomerization of O-Protonated (S)-trans-4-Hexen-3-ol. Journal of the American Chemical Society, 1997, 119, 4525-4534.	13.7	15
79	Gaseous protonated nitrosamide. A G2 theoretical study on the structure, stability, and interconversion of (H2Nî—¸NO) H+ isomers. Chemical Physics Letters, 1997, 267, 98-104.	2.6	17
80	Methylated NF3. A G2MS theoretical study on the structure, stability, and interconversion of the CH3â€“NF3+ and CH3Fâ€“NF2+ isomers. Chemical Physics Letters, 1997, 281, 431-437.	2.6	20
81	Aktivierung von Kohlendioxid durch Koordination mit Kationen in der Gasphase: SiF₃⁺â€vermittelte Carbonylierung von Arenen mit Kohlendioxid. Angewandte Chemie, 1996, 108, 2674-2676.	2.0	0
82	Protonated thiohypofluorous acid, FSH2+. Theoretically predicted to be stable and experimentally observed in the gas phase. Chemical Physics Letters, 1996, 253, 189-195.	2.6	10
83	Protonated methyl nitrite. A theoretical investigation on the structure and stability of (MeOî—¸NO)H+ and the proton affinity of ROî—¸NO (R = H, Me). Chemical Physics Letters, 1996, 258, 123-128.	2.6	15
84	Ionic Fluorination of Carbon Monoxide as a Route to Gasphase Carbonylation of Inert Cï£¿H and Nï£¿H Bonds. Chemistry - A European Journal, 1996, 2, 495-501.	3.3	35
85	Activation of Carbon Dioxide by Coordination with Cations in the Gas Phase: SiF3+-Mediated Coupling of CO2 and Aromatic Cï£¿H Bonds. Angewandte Chemie International Edition in English, 1996, 35, 2522-2524.	4.4	9
86	Structure and Stability of Isomeric C2GeH7+ Ions. An ab Initio Post-SCF Study. The Journal of Physical Chemistry, 1995, 99, 17724-17728.	2.9	8
87	Gaseous Fluorodiazonium Ions. Experimental and Theoretical Study on Formation and Structure of FN2+. Inorganic Chemistry, 1995, 34, 1325-1332.	4.0	25
88	The gaseous trifluorosilylxenon cation, F3SiXe+: a stable species with a siliconâ€“xenon bond. Journal of the Chemical Society Chemical Communications, 1995, , 773-774.	2.0	29
89	Gaseous Protonated Nitrosyl Fluoride. Experimental and Theoretical Characterization of Two Distinguishable Isomers, HONF+ and ONFH+, and Evaluation of the Barrier for Their Interconversion. The Journal of Physical Chemistry, 1994, 98, 2713-2718.	2.9	28
90	Gaseous F2NO+ Cations from the Addition of NF2+ to N2O. Structure and Mechanism of Formation. The Journal of Physical Chemistry, 1994, 98, 8009-8013.	2.9	22

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91	Uncommon electronic effects on the gas-phase BrÃ¸nsted acidity of isomeric hydroxyphenylium ions. Chemical Physics Letters, 1994, 229, 581-586.	2.6	3
92	An Extraordinarily Violent Molecular Dissociation: The Unprecedented Kinetic Energy Release in the Decomposition of HONF+, a Singly Charged Metastable Ion. Angewandte Chemie International Edition in English, 1994, 33, 123-125.	4.4	24
93	Eine auÃŸergewÃ¶hnlich heftige molekulare Dissoziation: beispiellose Freisetzung kinetischer Energie beim Zerfall von HONF, einem einfach geladenen, metastabilen Ion. Angewandte Chemie, 1994, 106, 104-106.	2.0	5
94	The addition of NF+2 to H2O as a route to gaseous protonated F2NOH. International Journal of Mass Spectrometry and Ion Processes, 1994, 130, 117-125.	1.8	17
95	Ionic Lewis superacids in the gas phase. Part 4. CF+3 initiated ion/molecule reaction patterns in the Î³-radiolysis of CF4/n-bases gaseous mixtures. International Journal of Mass Spectrometry and Ion Processes, 1994, 130, 207-222.	1.8	7
96	Experimental observation of stable cyanodiazonium ions, NCâ€“N2+. Journal of the Chemical Society Chemical Communications, 1994, , 2173-2174.	2.0	13
97	Gas-phase protonation of simple inorganic molecules: A stimulating interplay between theory and experiment. Organic Mass Spectrometry, 1993, 28, 1504-1511.	1.3	4
98	The NF2H+. and NH2F+. radical cations: conventional structures or ion-molecule complexes? A GAUSSIAN-1 study. Chemical Physics Letters, 1993, 204, 53-58.	2.6	6
99	Ionic Lewis superacids in the gas phase. Part 1. Ionic intermediates from the attack of gaseous SiF+3 on n-bases. International Journal of Mass Spectrometry and Ion Processes, 1993, 124, 21-36.	1.8	29
100	Ionic Lewis superacids in the gas phase. Part 2. Reactions of gaseous CF+3 with oxygen bases. International Journal of Mass Spectrometry and Ion Processes, 1993, 127, 123-135.	1.8	20
101	Ionic Lewis superacids in the gas phase. Part 3. Reactions of gaseous CF+3 with nitrogen bases. International Journal of Mass Spectrometry and Ion Processes, 1993, 127, 137-146.	1.8	15
102	Gas-phase protonation of spiropentane. A novel entry into the C5H9+ potential energy surface. Journal of the American Chemical Society, 1993, 115, 10338-10347.	13.7	14
103	Gas-phase ion chemistry of nitramide. A mass spectrometric and ab initio study of nitramide (H2N-NO2) and the H2N-NO2.+, [H2N-NO2]H+, and [HN-NO2]- ions. Journal of the American Chemical Society, 1993, 115, 12398-12404.	13.7	22
104	Concerning the proton affinity of hydrazoic acid and methyl nitrate. Journal of Organic Chemistry, 1993, 58, 3639-3642.	3.2	4
105	Gas-phase ion chemistry of cyanamide. A mass spectrometric and ab initio study of gaseous [H2N-CN].+, [H2N-CN]H+, and [HN-CN]- ions. The Journal of Physical Chemistry, 1993, 97, 4239-4245.	2.9	24
106	Structure and stability of various (C2,H5,Ge)+ ions: an ab initio molecular orbital study. The Journal of Physical Chemistry, 1993, 97, 4945-4950.	2.9	9
107	Structure and stability of H4NO4+ ions: an ab initio theoretical investigation. The Journal of Physical Chemistry, 1992, 96, 4354-4358.	2.9	8
108	Experimental and ab initio MO studies on [H2,N,O]+ ions in the gas phase: characterization of the isomers H2NO+, HNOH+ and NOH2+ and the mechanism of unimolecular dehydrogenation of [H2,N,O]+. The Journal of Physical Chemistry, 1992, 96, 4841-4845.	2.9	21

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109	Gas-phase protonation of nitrosyl hydride: a GAUSSIAN-1 ab initio MO study of the structure, stability, and unimolecular interconversion processes of various [H2,N,O]+ isomers. The Journal of Physical Chemistry, 1992, 96, 2100-2103.	2.9	12
110	Nitrogen versus fluorine protonation of nitrogen fluoride in the gas-phase. A combined mass spectrometric and Gaussian-1 ab initio MO study reveals the existence of two distinct isomers F3NH+ and F2N-FH+. Journal of the American Chemical Society, 1992, 114, 2806-2810.	13.7	57
111	Positive ion chemistry of gaseous boric and polyboric acids. International Journal of Mass Spectrometry and Ion Processes, 1992, 117, 47-63.	1.8	7
112	Gas-phase ion chemistry of H3BO3. Protonated orthoboric, metaboric and polyboric acids, and their anions in the gas phase. Journal of the Chemical Society Chemical Communications, 1991, , 66-68.	2.0	9
113	Gas-phase heteroaromatic substitution. 13. A quantitative application of the curve-crossing reactivity model to heteroaromatic substitution. Journal of the American Chemical Society, 1991, 113, 4550-4557.	13.7	11
114	H2NO2+ ions in the gas phase: a mass spectrometric and post-SCF ab initio study. The Journal of Physical Chemistry, 1991, 95, 9782-9787.	2.9	32
115	Bonding in square-planar MCl(CX)[P(i-Pr)3]2 complexes of rhodium and iridium (X = O and CH2) studied by UV photoelectron spectroscopy and DV-XÎ± calculations. Journal of Organometallic Chemistry, 1990, 382, 445-454.	1.8	4
116	Gas-phase heteroaromatic substitution. 8. Electrophilic attack of ethyl cation on pyrrole, N-methylpyrrole, furan, and thiophene. Journal of the American Chemical Society, 1990, 112, 3064-3068.	13.7	10
117	Relative stability of isomeric methyl nitrate cations (CH3NO3)H+. Journal of the Chemical Society Perkin Transactions II, 1989, , 413.	0.9	7
118	Tin-sulfur and tin-selenium bonding in some tin(IV) compounds studied by UV photoelectron and NMR spectroscopy and pseudopotential ab initio calculations. Organometallics, 1988, 7, 262-266.	2.3	5
119	A comparative study of gas phase aromatic desilylation and detertbutylation by charged electrophiles. Canadian Journal of Chemistry, 1988, 66, 3099-3107.	1.1	37
120	Ring-size effects on the ionization potentials of N-substituted azacycloalkanes. Journal of the Chemical Society Perkin Transactions II, 1986, , 667.	0.9	14
121	Evidence for Ïƒ-Ï€ interaction in some allyltin compounds. UV photoelectron spectroscopy and LCBO calculations. Journal of Organometallic Chemistry, 1986, 315, 287-297.	1.8	10